

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012717\
 Data File : BF092594.D
 Acq On : 27 Jan 2017 23:43
 Operator : SJ/MA
 Sample : I1448-08
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHA-BL-01-0-2

Quant Time: Jan 28 04:38:33 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 24 13:48:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	179963	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	695636	20.00	ng	0.00
38) Acenaphthene-d10	10.03	164	331415	20.00	ng	0.00
63) Phenanthrene-d10	11.53	188	446552	20.00	ng	0.00
75) Chrysene-d12	14.17	240	359775	20.00	ng	0.00
86) Perylene-d12	15.67	264	268785	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1468162	126.93	ng	0.05
7) Phenol-d6	6.62	99	1812475	123.02	ng	0.03
23) Nitrobenzene-d5	7.55	82	1324547	104.01	ng	0.01
41) 2,4,6-Tribromophenol	10.83	330	305946	135.47	ng	0.01
44) 2-Fluorobiphenyl	9.36	172	2022664	112.86	ng	0.00
78) Terphenyl-d14	13.12	244	1308674	96.86	ng	0.00
Target Compounds						
10) Phenol	6.64	94	135314	7.61	ng	Qvalue # 74
49) Dimethylphthalate	9.74	163	351825	16.69	ng	100
70) Phenanthrene	11.55	178	112059	4.56	ng	96
74) Fluoranthene	12.74	202	246143	10.23	ng	95
77) Pyrene	12.97	202	207581	7.96	ng	98
80) Benzo(a)anthracene	14.16	228	71001	3.67	ng	# 93
82) Chrysene	14.19	228	89516	4.34	ng	# 96
87) Benzo(b)fluoranthene	15.22	252	87396m	5.06	ng	
89) Benzo(a)pyrene	15.60	252	45664	3.13	ng	# 67
91) Benzo(g,h,i)perylene	17.62	276	28702	2.40	ng	# 66

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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